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Tommy Dessup, Christophe Coste, Michel Saint-Jean. Hysteretic and intermittent regimes in the subcritical bifurcation of a quasi 1D system of interacting particles. 2015. hal-01216045v1

HAL Id: hal-01216045 https://hal.science/hal-01216045v1

Preprint submitted on 16 Oct 2015 (v1), last revised 29 Nov 2016 (v2)

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Hysteretic and intermittent regimes in the subcritical bifurcation of a quasi 1D system of interacting particles

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Abstract

In this article, we study the effects of a white gaussian additive thermal noise on a subcritical pitchfork bifurcation. We consider a quasi-1D system of particles that are transversally confined, with short range (non Coulombic) interactions and periodic boundary conditions in the longitudinal direction. In such systems, there is a structural transition from a linear order to a staggered raw, called the zigzag transition. There is a finite range of transverse confinement stiffnesses for which the stable configuration at zero temperature is a localized zigzag pattern surrounded by aligned particles, which evidences the subcriticality of the bifurcation. We show that these configurations remains stable for a wide temperature range. At zero temperature, the transition between a straight line and such localized zigzag patterns is hysteretic. We have studied the influence of the thermal noise on the hysteresis loop. Its description is more difficult than at T = 0 K since thermally activated jumps between the two configurations always occur and the system can never stay forever in a unique metastable state. Two different regimes have to be considered according to the temperature value with respect to a critical temperature $T_c(\tau_{obs})$ that depends on the observation time τ_{obs} . An hysteresis loop is still observed at low temperature, with a width that decreases as the temperature increases toward $T_c(\tau_{obs})$. In contrast for $T > T_c(\tau_{obs})$ the memory of the initial condition is lost by stochastic jumps between the configurations. The study of the mean residence times in each configurations gives a unique opportunity to precisely determine the barrier height that separates the two configurations, without knowing the complete energy landscape of this many-body system. We also show how to reconstruct the hysteresis loop which would exist at T = 0 K from high temperature simulations.

PACS :

05. Statistical physics, thermodynamics, and nonlinear dynamical systems
63.70.+h Statistical mechanics of lattice vibrations and displacive phase transitions
05.40.-a Fluctuation phenomena, random processes, noise, and Brownian motion

I. INTRODUCTION

There is much recent interest in physical systems constituted by interacting particles confined in narrow channels, in quasi-one-dimensional geometry. Their typical size extends on several orders of magnitude, from laser-cooled ions confined in Paul traps [1–14] and polar molecules [15] to optically confined paramagnetic colloidal particles [16], to plasma dusts in electrostatic traps [17–21], and to electrostatically interacting macroscopic beads [14, 22–24]. The transverse confining potential is well described as harmonic (see [6] for an anharmonic potential), while the repulsive interactions may be either described by the Coulomb potential [1–13], a dipolar interaction [15, 16], a Yukawa potential $U_0 \exp(-r/\lambda_0)/r$ for screened electrostatic interaction [17–21] or interaction potential $U_0 K_0(r/\lambda_0)$ [14, 22–24], where $K_0(.)$ is the modified Bessel function of order 0, and where U_0 and λ_0 are respectively the intensity and the range of interaction.

The behavior of these systems results from the interplay between the repulsive inter particle interactions and the potential that confines the particles in the transverse direction of the trap, which may be characterized by its stiffness β . When β is large enough, all particles are aligned along the trap axis at equilibrium. As β decreases below a critical stiffness β_{ZZ} , the equilibrium pattern undergoes a structural phase transition. The characteristics of the transition depend on the range of the inter particles interactions, on the boundary conditions in the longitudinal direction and on the system size.

For an infinite interaction range $(\lambda_0 \to \infty)$, typically for Coulombic interactions, a system always evolves from a straight line toward one of two regular symmetric staggered rows. This transition is therefore called the *zigzag transition*. Its relevant order parameter is the height h defined as the distance between any particle and the longitudinal axis. A suitable control parameter is the dimensionless transverse stiffness $\epsilon = 1 - \beta/\beta_{ZZ}$ (see also [25]). The bifurcation diagram $h(\epsilon)$ corresponds to a supercritical pitchfork bifurcation : The zigzag phase spontaneously breaks the mirror symmetry, the zigzag phase is stable for $\epsilon \ge 0$, and its amplitude $h(\epsilon)$ is proportional to $\epsilon^{1/2}$ [1–4, 13, 14] [see Fig. 1-(a)].

For longitudinally confined systems and whatever the inter particle interaction range, the same kind of transition can be observed. Because of the longitudinal boundary conditions, the two symmetric zigzag configurations are now inhomogeneous, with the highest distance h_m always localized at the center of the finite cell, and a zigzag amplitude that symmetrically

decreases toward both edges of the channel. In this case, the relevant order parameter is the maximum zigzag amplitude h_m . Its evolution with the dimensionless transverse stiffness ϵ evidences the supercritical nature of the bifurcation [3, 4, 7, 20, 22, 26].



FIG. 1: (Color online) (a) Supercritical pitchfork bifurcation diagram at T = 0 K. (b) Subcritical pitchfork bifurcation diagram at T = 0 K. The control parameter $\epsilon = 1 - \beta/\beta_{ZZ}$ is the dimensionless transverse confinement stiffness, which increases from left to right. For the zigzag transition the order parameter h_m is the maximum height of the equilibrium configuration. The red (dark grey) continuous and dashed lines correspond respectively to stable and unstable configurations. The cyan (light grey) line displays an hysteresis loop.

In a finite periodic system with N particles contained in a cell of length L, when the inter particles interaction is of finite range ($\lambda_0 < \infty$), the zigzag transition is strongly modified. This is relevant for experiments in annular cells [19, 27] and for simulations [14, 24]. In these systems, when $\lambda_0/d \gg 1$ where $d \equiv L/N$ is the mean inter particles distance, the zigzag transition is still a supercritical pitchfork bifurcation (we assume N even to allow for an homogeneous zigzag phase in the cell). For smaller values of λ_0/d , the description of the zigzag transition is more complex. As the confinement decreases, the system first evolves from a linear configuration to an homogeneous zigzag. This happens for $\beta = \beta_{ZZ}$ $(\epsilon = 0)$. Our previous linear stability analysis allows to identify the domain of stability of this homogeneous zigzag configuration. This domain extends from $\epsilon = 0$ up to a positive $\epsilon > 0$ that depends on the particles number N at fixed λ_0/d . It is a strongly decreasing function of N and vanishes in the thermodynamic limit $(N \to \infty \text{ at fixed } d)$ [14, 24]. Then for smaller confinement, an inhomogeneous zigzag pattern is observed in which particles in a line coexist with particles organized in a distorted zigzag (see Fig. 2). We have called this pattern a bubble [19, 24]. Eventually, at very small confinement ($\epsilon \leq 1$) an homogeneous zigzag is restored.

Since for inhomogeneous patterns the maximum distance h_m can be taken as an order parameter, we have performed numerical simulations to determine the variations of $h_m(\epsilon)$ for increasing and decreasing ϵ . In Ref. [24], we have provided evidences of an hysteretic behavior : The transition from a line toward a bubble, under decreasing confinement, takes place at a smaller confinement than the transition from a bubble toward a line, under increasing confinement. These observations of inhomogeneous equilibrium patterns and of an hysteresis loop all indicate a transition described by a *subcritical* pitchfork bifurcation [see Fig. 1 (b)] for finite range interactions and periodic boundary conditions in the longitudinal direction.

From the theoretical point of view, by considering the vibrational modes of homogeneous patterns in the case of an infinite system of particles, we have shown that $\omega(q = \pi/d)$, the frequency associated to the alternate transverse displacement of neighboring particles, vanishes at the confinement threshold. This means that at the transition the displacements associated to this soft-mode are quenched, resulting in the homogeneous zigzag pattern [14, 22]. For an infinite system, this soft mode coexists with the Goldstone mode associated to the translational invariance of the infinite line. Extending the analysis to small nonlinearities, in the thermodynamic limit and near the bifurcation threshold [24], we have exhibited the key role of the nonlinear coupling between the soft mode at the bifurcation and the Goldstone mode linked to the translational invariance. We have obtained the normal form of the bifurcation, and the conditions required to observe a supercritical or a subcritical pitchfork bifurcation.

In periodic systems there is also a Goldstone mode linked to the rotational invariance of the periodic cell. We have established that for the parameters used in our experiments and in our simulations the bifurcation is always subcritical. A quantitative description of the nonlinear localized zigzag patterns (bubbles) observed in the simulations is deduced from the explicit normal form, in excellent agreement with the simulations (see Fig. 2, the Appendix B and Figs. (10)-(12) of Ref. [24]).

The discussion so far concerns a structural transition observed at temperature T = 0 K. Beyond this first approach an interesting question is to study the validity and the limits of this description for systems submitted to a thermal noise. In this paper the focus is set on this structural transition at finite temperature. The bubble diffusion induced by thermal activation will be presented in a forthcoming paper [28].

In the literature there are some theoretical studies devoted to the thermal effects on supercritical bifurcations [29, 30]. Their results are in good agreement with our numerical study of the thermal noise effects on the zigzag transition for longitudinally confined systems [23]. More recently, this approach has been extended to subcritical bifurcations. In particular, it was shown that temperature affects the hysteresis loop [31]. However the expressions developed in these study are essentially formal, and poorly confronted with actual data.

In this article, we study the effects of a white gaussian additive thermal noise on a subcritical pitchfork bifurcation. The temperature range that we consider is selected according to two limiting constraints. At too low temperature no effects can be expected since the particles only vibrate around their equilibrium position associated to the configurations obtained at T = 0 K. On the other hand, at too large temperature, the amplitude of transverse fluctuations are of the same order as the bubble height so that the instantaneous patterns are completely random.

In section II we first discuss the stability of the bubbles with increasing thermal noise, and the relevance of our theoretical description of the bubbles to systems at finite (nonzero) temperature. Then we consider the hysteresis loop between a straight line and a bubble observed at very low temperature [24]. At T = 0 K the system is in one of the two metastable configurations according to its history : The initial configuration and the direction of change of the control parameter ϵ . It's not the case anymore at $T \neq 0$. Indeed inside the hysteresis loop the temperature always induces thermal jumps between the two configurations. According to the Kramers theory [32], these jumps are characterized by two transition times $\tau_{L-B}(\epsilon, T)$ (from line to bubble) and $\tau_{B-L}(\epsilon, T)$ (from bubble to line). The existence of these times suggests the important role played on the bifurcation description by the observation time τ_{obs} during which ϵ is constant. As long as τ_{obs} is smaller than the characteristic transition times, the jumps between the configurations are not observed. In contrast the jumps have to be taken into account once one of these Kramers' times is smaller than τ_{obs} . By comparing the three characteristic times τ_{obs} , $\tau_{L-B}(\epsilon, T)$ and $\tau_{B-L}(\epsilon, T)$, we may define two distinct regimes : The *hysteretic regime* for which an hysteresis loop may still be defined, and the *intermittent regime* for which the thermally activated jumps blur the hysteresis loop. Using the probability density function (PDF) of the instantaneous bubble amplitude $h_m(\epsilon, t)$, we are able to obtain an efficient description of the bifurcation at $T \neq 0$. In particular, this allows a detailed study of the relation between τ_{obs} and the hysteretic and intermittent regimes. The evolution of the hysteresis loop with temperature and the intermittency effects will be exhibited.

In Sec. III we focus on a temperature range that is low enough for the hysteresis thresholds to be defined. We show that the hysteresis loop width decreases as the temperature increases. A threshold temperature $T_c(\tau_{obs})$ beyond which no hysteresis can be observed is determined as a function of τ_{obs} . The rough features of this function are recovered with the help of a qualitative analysis. This critical temperature is linked to a particular value ϵ_M of the confinement, which is defined as a Maxwell point for the relevant energy landscape and does not depend on τ_{obs} . A numerical value of ϵ_M is provided.

In Sect. IV, we focus on the intermittent regime observed at higher temperature. We show that in this regime a mean residence time for each of the configurations may be defined and measured on the simulations data. The variations of this residence times with temperature can be used to determine the main features of the system energy landscape. We propose although a procedure to reconstruct the hysteresis diagram at T = 0 K from measurements performed at $T \neq 0$. Lastly, we obtain two other numerical values for the confinement ϵ_M , with two independent methods, in very good quantitative agreement with the previous one.

The simulations results presented here have been obtained by molecular dynamics with small dissipation in order to emphasize the noise effects on the particle dynamics. The details about simulations are given in Appendix A. For the sake of completeness the theoretical description of the bubbles is summarized in Appendix B.

II. EVOLUTION OF BUBBLE PATTERNS WITH THERMAL NOISE

A. Bubbles at finite temperature

At T = 0 K, for each confinement ϵ corresponding to a bubble state, the particles are organized according to a unique stationary equilibrium pattern. Their shape is accurately deduced from the normal form (B1) given in Appendix B. The theoretical bubbles shape is given in Eqn. (B3) and the determination of an *effective bifurcation parameter*, mandatory for finite systems with periodic longitudinal boundary conditions, is made from Eqns. (B5) and (B6), thus without any fitting parameter (see also Ref. [24]). The first question that arises is about the possible persistence of these patterns at higher temperatures and correspondingly the relevance of the theoretical analysis to their description.

In the Figs. 2 (a)–(c), we show instantaneous pictures of the periodic cell for a constant transverse confinement and increasing temperatures, which are at least two orders of magnitude higher than in our previous study [24]. In each case a bubble is evidenced, and whatever the temperature its shape compares very accurately with the unique theoretical shape deduced from the known value of ϵ . This proves that at these temperatures these coherent patterns are not destroyed by the thermal fluctuations. To emphasize this point, we plot in Figs 2 (d)–(f) the superposition of all the positions reached by particles during a typical run. At $T = 10^9$ K and $T = 10^{10}$ K [Figs 2 (d)–(e)], the bubble shape is basically preserved during the whole simulation run. The bubbles are actually stable equilibrium patterns, the thermal fluctuations only inducing random motions of the particles around their equilibrium positions, with an amplitude that increases with temperature. At higher temperature $[T = 10^{11}$ K, see Figs 2 (f)] the bubble pattern remains recognizable even if it is almost completely blurred because of the fluctuations.

B. Hysteretic and intermittent regimes

The systems presenting a subcritical pitchfork bifurcation exhibit a finite range of ϵ for which the configuration may be a line or a bubble (see Appendix B for a brief summary). These metastable configurations are associated to the energy minima $E_L(\epsilon)$ for the line and $E_B(\epsilon)$ for the bubble. A thermal fluctuation may activate a jump from one configuration to the other at the expense of a maximum of energy $E_M(\epsilon)$. We insist on the fact that this maximum is not provided by our continuous model, since the knowledge of the full energy landscape require the complete solution of the discrete N-body problem. We discuss this point in details in Sec. IV.

Let us first consider the system at T = 0 K, and let us look at the configuration evolution when ϵ increases for a system that is initially in the linear configuration. The system remains in this configuration as long as $E_L(\epsilon)$ is a local minimum. When ϵ reaches the value $\epsilon_{up}(T = 0)$, the minimum $E_L(\epsilon_{up})$ and the maximum $E_M(\epsilon_{up})$ disappear, and the



FIG. 2: (Color online) Particles position (x_i, y_i) (mm) of 128 particles in the bubble configuration ($\epsilon = 0.02$ and d = 1.875 mm), for three temperatures $T = 10^9$, 10^{10} and 10^{11} K from left to right. In the top plots the green dots correspond to an instantaneous particles position and the dashed red line is the theoretical bubble shape. In the bottom plots the blue dots (dark gray) are the mean values of the positions, the light gray dots corresponds to superposition of all instantaneous positions recorded during the simulation run and the solid red line is the theoretical bubble shape.

system jumps to a bubble configuration of smaller energy $E_B(\epsilon_{up})$. In the same manner, by considering a system initially in the bubble configuration and a decreasing ϵ we can define $\epsilon_{down}(T=0)$. These two thresholds characterize the width $\epsilon_{up}(T=0) - \epsilon_{down}(T=0)$ of the hysteresis loop at T=0 K.

At $T \neq 0$ K, these jumps are thermally activated. The jump from a linear to a bubble configuration is characterized by a Kramers' activation time $\tau_{L-B}(\epsilon, T) \propto \exp(\Delta U_{L-B}(\epsilon)/k_B T)$ where $\Delta U_{L-B}(\epsilon) = E_M(\epsilon) - E_L(\epsilon)$, whereas the jump from a bubble to a linear configuration is characterized by $\tau_{B-L}(\epsilon, T) \propto \exp(\Delta U_{B-L}(\epsilon)/k_B T)$ where $\Delta U_{B-L}(\epsilon) = E_M(\epsilon) - E_B(\epsilon)$. Thus the probability of a jump from one configuration to the other one is a function of both ϵ and T.

Because of the existence of these Kramers' times, the hysteresis description depends on the observation time τ_{obs} . For an infinite τ_{obs} , whatever ϵ , the Kramers' times are necessarily smaller than τ_{obs} and both configurations are observed, independently of the initial state. So we cannot define the transition thresholds as for T = 0 K and the hysteresis loop is meaningless. However, in practice any actual numerical or experimental observation implies a finite τ_{obs} . The hysteresis thresholds have to be redefined and depend on this observation time. A threshold can be defined if a jump from an initial configuration to the other one occurs during the observation time, while the reverse jump does not occur because the corresponding Kramers' time is too long with respect to τ_{obs} . In this case, speaking of an hysteresis loop that depends on the temperature makes sense. In contrast, when the characteristic times for both jumps are smaller than τ_{obs} , the two configurations are observed during τ_{obs} , whatever the initial configuration of the system. The memory of the initial configuration is not conserved, so that no hysteresis loop can be defined. We will speak in this case of an *intermittent regime*.

C. Definition of an order parameter in presence of thermal noise

In order to describe the transition it is mandatory to define the order parameter $h_m(\epsilon)$ for each confinement ϵ , that characterizes accurately and unambiguously the configurations explored by the system. The definition of this parameter at finite temperature is not obvious and requires an adapted procedure.

The simulations only provide instantaneous configurations. Such configurations are plotted for three different temperatures in Fig. 2 (a)–(c). They are very sensitive to the thermal fluctuations, however they are all in close agreement with the same theoretical description, independently of the temperature, as shown in Sec. II A. Thus, the main question is how to determine an amplitude $h_m(\epsilon)$ which could well characterize this configuration. We could be tempted to define $h_m(\epsilon)$ as the maximum distance between the particles and the longitudinal axis, measured from time averaged configurations. This procedure makes sense at low temperature, but it introduces important errors at high temperature. Indeed, at small temperature, we see that the averaged configurations [see Fig. 2 (d) and (e)] are very close to the instantaneous ones [see Fig. 2 (a) and (b)]. In contrast the averaged configuration at the highest temperature, shown in Fig. 2 (f), has nothing in common with the instantaneous configuration of Fig. 2 (c). Thus the strong discrepancy between the averaged configuration of Fig. 2 (f) and those of Fig. 2 (d) and (e) is a spurious effect which prevents the definition of $h_m(\epsilon)$ from the averaged configuration.

We have therefore developed a consistent process to determine the amplitude $h_m(\epsilon)$ at every temperature. In order to take advantage of the improvement in the statistics implied by the duration of the whole simulation, while avoiding averaging artifacts, we define $h_m(\epsilon, t)$ as the maximum distance between the particles and the cell axis at time t, and we consider its probability density function (PDF). The PDFs are shown in Fig. 3 for two temperatures and increasing values of ϵ , near the bifurcation threshold.

The position of the maximum of this distribution may be used as an order parameter. Let us first consider a strong confinement ϵ for which the particles are along a straight line at zero temperature, so that $h_m(\epsilon, T = 0, t) = 0$ and the PDF is the Dirac δ -function $\delta(h)$. At finite temperature, the maximum height $h_m(\epsilon, T > 0, t)$ is strictly positive at each time and increases with temperature [see Fig. 3 (a)]. Thus the corresponding PDF is slightly shifted above zero and is broadened around the most probable value $h_m(\epsilon, T > 0)$. Note that from now on we emphasize the dependence of h_m upon the temperature.

Then, for a confinement such that the stable pattern at T = 0 is a bubble, the PDF is a Dirac δ -function centered on a strictly positive value that is the bubble amplitude. At finite temperature, the PDF evidences a small shift toward higher values and a broadening, as was the case for the straight line [see Fig. 3 (c)]. The configuration is nevertheless easy to distinguish from a straight line, because despite the thermal fluctuations the most probable value $h_m(\epsilon, T > 0)$ is much higher than in the previous case.

Lastly, let us consider a confinement ϵ in the hysteretic domain at T = 0 K. At zero temperature, the distribution of $h_m(\epsilon, T = 0, t)$ corresponds to Dirac δ -functions centered around $h_m = 0$ if the initial configuration was a straight line or around $h_m > 0$ if the initial configuration was a bubble configuration. The PDF of $h_m(\epsilon, T, t)$ at finite temperature is shown in Fig. 3 (b). At low temperature ($T = 10^{10}$ K, cyan–light grey symbols) for which no jump between the two allowed configurations may be expected, the system simply keeps a memory of its initial configuration and explores the configuration space around only one stable attractor. Depending upon the initial condition, the PDF is a finite width distribution centered around a small value of h_m for the fluctuating straight line [Fig. 3 (b), cyan circles] or around a much larger value of h_m for the fluctuating bubble [Fig. 3 (b), cyan crosses]. At higher temperature, ($T = 10^{11}$ K, red–dark gray symbols) there are thermally activated jumps between the line and the bubble. This is the intermittent regime, for which the memory of its initial condition is lost by the system. The PDF evidences a camelback shape [see Fig. 3 (b), red–dark grey symbols]. The first peak corresponds to the fluctuating straight line configurations and the second peak corresponds to the fluctuating bubble configurations. Both peaks are broadened because of the thermal fluctuations, and the probability to observe an intermediate configuration that cannot be described neither as a line nor as a bubble is strictly positive.



FIG. 3: (Color online) PDF of the height $h_m(\epsilon, T, t)$ (in mm) for a system of 128 particles, for three different confinements : (a) $\epsilon = -0.05$; (b) $\epsilon = 0.001$; (c) $\epsilon = +0.05$. The red (dark grey) and cyan (light grey) dots correspond respectively to $T = 10^{10}$ K and $T = 10^{11}$ K. In (b) the open circles correspond to a case of line initial condition and the crosses to the case of bubble initial condition. In the inset of each plot, we show the corresponding point in the hysteresis loop measured at a very low temperature $T = 10^7$ K.

D. Evolution of the hysteresis with temperature

The thermal effects on the hysteresis diagram can be studied by following the evolution of the PDFs of $h_m(\epsilon, T, t)$ with the confinement ϵ for several temperatures. In Fig. 4 we display the reconstructed hysteresis diagram with the help of the PDFs, shown as density plots for each confinement ϵ .

For low temperature, the thermal energy is insufficient to induce intermittency during a step at constant ϵ of duration τ_{obs} . The hysteresis loop is still observed, and the diagram is only slightly modified with respect to the case of the very low temperature $T = 10^7$ K, as shown by Fig. 4 (a) and (b). The comparison between the hysteresis diagrams obtained at $T = 10^9$ K and $T = 10^{10}$ K (resp. Fig. 4-(a) and 4-(b)) shows that its width decreases as the temperature increases. This point will be discussed in details in the next section III.

For higher temperature, the intermittent behavior strongly modifies the transition diagram, as shown by Fig. 4 (c) and (d). Now the thermally activated jumps between the linear configuration and the bubble happen in the range $\epsilon_{down}(T=0) < \epsilon < \epsilon_{up}(T=0)$. At



FIG. 4: Hysteresis diagrams reconstructed from the PDF of $h_m(\epsilon, T, t)$, measured at temperatures $T = 10^9$ K (a), $T = 10^{10}$ K (b), $T = 10^{11}$ K (c) and $T = 2 \times 10^{11}$ K (d). The solid red line represents the hysteresis loop obtained at very low temperature $T = 10^7$ K. The PDF are displayed as density plots for each ϵ and correspond to the blue scales on the right. For $T = 10^9$ K and $T = 10^{10}$ K the system keeps the memory of its initial condition, and the evolution of ϵ during the whole simulation corresponds to the red arrows. At $T = 10^{11}$ K and $T = 2 \times 10^{11}$ K all memory is lost by the system, and the diagrams do not depend on the evolution of ϵ during the whole simulation.

 $T = 10^{11}$ K and $T = 2 \times 10^{11}$ K, during a step at constant ϵ of duration τ_{obs} , the system looses all memory of its initial state and the hysteresis loop is no more observed.

It is possible to get a qualitative picture of these hysteresis diagram. Whereas at T = 0 K the energy of the system is either $E_L(\epsilon)$ or $E_B(\epsilon)$, the thermal agitation results in an exploration of the neighborhoods of $E_L(\epsilon)$ and $E_B(\epsilon)$ by the energy of the system. This induces a broadening of the PDF of $h_m(\epsilon, T, t)$. As said before, we do not know the complete energy landscape. We may nevertheless obtain a rough estimate of the distribution of $h_m(\epsilon, T, t)$ by considering only homogeneous configurations. In this simpler model the expression of the energy reads :

$$E(\epsilon, h) = \frac{\epsilon \beta_{ZZ}}{2} h^2 + \frac{g_3}{4} h^4 - \frac{g_5}{6} h^6$$
(1)

where ϵ is the order parameter, h the observable (the amplitude of the homogeneous zigzag) and β_{ZZ} and the g_i are positive constants [See Eqn. (B1) of Appendix B]. The hysteresis loop at T = 0 is then characterized by the existence of two local energetic minima, one for h = 0 noted $E_0(\epsilon)$ and one for $h \neq 0$ noted $E_1(\epsilon)$. In order to exhibit the broadening we may estimate the extension of h reached by the system thanks to the thermal fluctuations, by numerically extracting h from the equations $E(\epsilon, h) = E_0(\epsilon) + k_B T$ near the solution h(T = 0) = 0 and $E(\epsilon, h) = E_1(\epsilon) + k_B T$ near the solution $h(T = 0) \neq 0$. The domain explored by h calculated using this procedure are plotted in Fig. 5 when an hysteresis loop is observed at low temperature [Fig. 5 (a)], and when an intermittent regime is observed at high temperature [Fig. 5 (b)]. They display the same qualitative features as those obtained in the simulations for inhomogeneous configurations (Fig. 4).

III. EFFECT OF THE THERMAL NOISE ON THE HYSTERESIS THRESHOLDS

We now look in details at the influence of the thermal noise on the hysteresis loop for a given observation time τ_{obs} during which ϵ is kept constant.

At T = 0 K the thresholds $\epsilon_{down}(T = 0)$ and $\epsilon_{up}(T = 0)$ only depend on the stability of the equilibrium configurations. At finite temperature, the thresholds $\epsilon_{down}(T, \tau_{obs})$ and $\epsilon_{up}(T, \tau_{obs})$ also depend on the temperature and on the observation time. Let us first consider a system initially in the linear configuration and increase ϵ by successive steps of duration τ_{obs} . For the *i*-th step, $\epsilon = \epsilon_i$, the system stays in a straight line configuration as long as the Kramers' transition time $\tau_{L-B}(\epsilon_i, T) > \tau_{obs}$, because its initial configuration at the step (i-1)is a line, and τ_{obs} is insufficient for the transition to happens. The transition toward the bubble configuration takes places when $\tau_{L-B}(\epsilon_i, T) \sim \tau_{obs}$ since the observation of a jump toward the bubble configuration during τ_{obs} is now more likely. Once the transition is achieved, if $\tau_{B-L}(\epsilon_i, T) > \tau_{obs}$ the system remains in the bubble configuration. Thus $\tau_{L-B}(\epsilon_{up}, T) \sim \tau_{obs}$ defines the hysteresis threshold $\epsilon_{up}(T, \tau_{obs})$ providing that $\tau_{B-L}(\epsilon_{up}, T) > \tau_{obs}$ at the temper-



FIG. 5: (Color online) Qualitative estimate of the thermal broadening for a subcritical pitchfork bifurcation. For each ϵ the red (grey) shaded segment quantify the reachable values of h. The solid black lines display the stable homogeneous states, the dotted black lines the unstable states and the black arrows the hysteresis loop at T = 0. (a) : At low temperature the two explored regions corresponding to the two equilibrium states are always distinct and the hysteresis loop remains as represented by the solid red line. This is to be compared to Figs. 4 (a) and (b). (b) : At high temperature the two explored area merge and the hysteresis loop is blurred. This is to be compared to Figs. 4 (c) and (d).

ature T. Note that since $\tau_{L-B}(\epsilon, T) \propto \exp(\Delta U_{L-B}(\epsilon)/k_B T)$ decreases with temperature, the condition $\tau_{L-B}(T, \epsilon_{up}) \sim \tau_{obs}$ implies that $\epsilon_{up}(T, \tau_{obs})$ is a decreasing function of the temperature. Similarly, following the evolution of an initial bubble configuration as ϵ decreases, we can define the threshold $\epsilon_{down}(T, \tau_{obs})$ (with $\tau_{B-L}(\epsilon_{down}, T) \sim \tau_{obs}$ and $\tau_{L-B}(\epsilon_{down}, T) > \tau_{obs}$). In the opposite way, this threshold increases with temperature.

The Fig. 6 represents $\epsilon_{down}(T, \tau_{obs})$ and $\epsilon_{up}(T, \tau_{obs})$ as functions of T, from our simulation data. The variations of $\epsilon_{down}(T, \tau_{obs})$ and $\epsilon_{up}(T, \tau_{obs})$ with the temperature agree with the above analysis. The width of the hysteresis loop, $\epsilon_{up}(T, \tau_{obs}) - \epsilon_{down}(T, \tau_{obs})$, eventually vanishes at a threshold temperature $T_c(\tau_{obs})$. This temperature is defined by $\tau_{L-B}(\epsilon_{down}, T_c) \sim \tau_{B-L}(\epsilon_{up}, T_c) \sim \tau_{obs}$.

A qualitative picture of the variation of T_c with τ_{obs} may be obtained. For an infinite observation time, intermittency is always present, thus no hysteresis loop can be defined at finite temperature so that $T_c(\tau_{obs} \to \infty) \to 0$. In order to analyze the variation of $T_c(\tau_{obs})$ for finite τ_{obs} we may derive the evolution of $\epsilon_{down}(T, \tau_{obs})$ and $\epsilon_{up}(T, \tau_{obs})$ from the two



FIG. 6: (Color online) Evolution of the thresholds $\epsilon_{up}(T, \tau_{obs})$ (blue dots) and $\epsilon_{down}(T, \tau_{obs})$ (red squares) with temperature.

equations :

$$k_B T \ln\left(\frac{\tau_{obs}}{\tau_0}\right) = \Delta U_{L-B}(\epsilon_{up}),$$
(2)

$$k_B T \ln\left(\frac{\tau_{obs}}{\tau_0}\right) = \Delta U_{B-L}(\epsilon_{down}),\tag{3}$$

where for the sake of simplicity the characteristic time τ_0 introduced by Kramers [32] may be considered constant.

Up to now these equations are only formal because we do not know the energy gaps. As before, we get a qualitative interpretation by considering homogeneous patterns. We show in Fig. 7 the evolution of $\epsilon_{up}(T, \tau_{obs})$ and $\epsilon_{down}(T, \tau_{obs})$ calculated for three ratios τ_{obs}/τ_0 . They are obtained by directly solving the equations (2) and (3) where the unknown energetic barriers ΔU_{L-B} and ΔU_{B-L} are estimated for homogeneous patterns from the simple energetic expression of Eqn. (1). We recover the qualitative features of the functions $\epsilon_{down}(T, \tau_{obs})$ and $\epsilon_{up}(T, \tau_{obs})$ measured from the simulations data, shown in Fig. 6. This is also consistent with simulations results presented by Agez et al. [31] see their Fig. 15.

The crossing of the plots $\epsilon_{down}(T, \tau_{obs})$ and $\epsilon_{up}(T, \tau_{obs})$ happens by construction for $T = T_c(\tau_{obs})$. It takes place at a confinement $\epsilon = \epsilon_M$ which does not depend on τ_{obs} . This particular value ϵ_M is defined as the confinement for which both energetic barriers are equals :

$$\Delta U_{L-B}(\epsilon_M) = \Delta U_{B-L}(\epsilon_M), \tag{4}$$

which suggest to call it the *Maxwell point*. Note that this equation and thus the value of ϵ_M are independent of the temperature. As shown in Fig. 6 the value of ϵ_M is 0 ± 0.01 .



FIG. 7: (Color online) Theoretical variation of the thresholds $\epsilon_{up}(T, \tau_{obs})$ (cyan, light grey) and $\epsilon_{down}(T, \tau_{obs})$ (red, dark grey) with the temperature, for observation times $\tau_{obs}/\tau_0 = 0.5$ (dashed line), $\tau_{obs}/\tau_0 = 1$ (solid line) and $\tau_{obs}/\tau_0 = 4$ (dotted line). The critical temperature $T_c(\tau_{obs})$ decreases as τ_{obs} increases. The confinement ϵ_M corresponds to the Maxwell point at which both stable states have the same energy.

IV. INTERMITTENT REGIME AND ENERGY LANDSCAPE

Let us now consider the intermittent regime. Despite the large fluctuations induced by the thermal agitation, we have taken advantage of the unique opportunity given by the numerous transitions between configurations to determine the main features of the energy landscape (Maxwell point and energetic barrier height).

A. Residence time and energy landscape

For a temperature higher than $T_c(\tau_{obs})$ the intermittent regime is reached as soon as $\tau_{B-L}(\epsilon, T) < \tau_{obs}$ and $\tau_{L-B}(\epsilon, T) < \tau_{obs}$. An example of such intermittency is shown in Fig. 8 which presents a typical time evolution of the maximum height $h_m(\epsilon, T, t)$. The system stays for a finite time in one configuration and jumps quickly to the other configuration. Let τ_L be the time spent by the system in the straight line configuration, and τ_B the time spent by the system in the straight line configuration, are widely distributed, as shown in Fig. 9 (a) and (b).

From such distributions the mean residence times $\langle \tau_L \rangle$ and $\langle \tau_B \rangle$ may be calculated [33].



FIG. 8: (Color online) Typical evolution of $h_m(\epsilon, T, t)$ (in mm) with time (in s) for $\epsilon = 0.005$ with $T = 8.10^{10}$ K. The solid black lines indicate h_m values associated to the linear and bubble configurations for T = 0 K, the dashed black lines indicate the time average values $h_m(\epsilon, T)$ for the linear and bubble configurations at $T = 8.10^{10}$ K.

In order to check the relevance of these two times, the PDF obtained from simulations have to be compared to the homogeneous Poisson probability density $P(\tau) = e^{-\tau/\langle \tau \rangle}/\langle \tau \rangle$, which describes the probability distribution of residence times that are random and of small probability [32]. In Fig. 9 (a) and (b) we show that the PDF measured from the simulations is very well fitted by the expected probability distribution.

The knowledge of these mean residence times provides a lot of useful information. Obviously, it gives the most likely state of the system, which cannot be deduced from the transition diagram blurred at this high temperature [see, e.g. Fig. 4 (c) and (d)]. Moreover, the evolution of these residence times with the confinement ϵ gives another method to determine the Maxwell point. At the Maxwell point, we expect that $\langle \tau_L \rangle = \langle \tau_B \rangle$. We display in Fig. 9 (c) the evolution of the mean residence times $\langle \tau_L \rangle$ and $\langle \tau_B \rangle$ as a function of ϵ at a temperature $T = 8.10^{10}$ K. The two mean residence times are equal for $\epsilon_M = 0.003 \pm 0.004$, which is in very good agreement with our previous estimate deduced from the zero width of the hysteretic loop (see Fig. 6).

Furthermore, the energy barriers $\Delta U_{L-B}(\epsilon)$ and $\Delta U_{B-L}(\epsilon)$ for a given ϵ may be estimated from the evolution of residence times with temperature. The first step is to determine the evolution of the mean residence times with the temperature, at fixed ϵ , Fig. 10. The plots measured on the simulations data are very well fitted by a Kramers' law [34], which gives a reliable estimate for the energy barriers $\Delta U_{L-B}(\epsilon)$ and $\Delta U_{B-L}(\epsilon)$ for a given ϵ .



FIG. 9: (Color online) (a) Distribution of the residence times for a line. (b) Distribution of the residence times for a bubble. In both cases the confinement is $\epsilon = 0.0$ and $T = 8.10^{10}$ K. The red dashed line is a fit with the expected Poisson distribution, $P(\tau) = e^{-\tau/\langle \tau \rangle}/\langle \tau \rangle$ from which we deduce the relevant mean residence time $\langle \tau \rangle$. (c) Plots of the mean residence times $\langle \tau_L \rangle$ (blue dots) and $\langle \tau_B \rangle$ (red squares) as a function of ϵ at $T = 8.10^{10}$ K. The crossing of these plots defines the Maxwell point ϵ_M (dotted line).

When we repeat this process for several values of ϵ , we are able to plot the evolution of the energy barriers with the confinement ϵ , as shown in Fig. 11. As expected, these barriers are monotonic functions of ϵ , with $\Delta U_{L-B}(\epsilon)$ decreasing and $\Delta U_{B-L}(\epsilon)$ increasing. Note that these energetic barriers are independent of the temperature and observation time, as expected for these intrinsic parameters of the system. The crossing point of this two curves, for which the two barriers are equal, characterizes by definition the Maxwell point ϵ_M [see Eqn. (4)]. Its value measured on Fig. 11 is $\epsilon_M = 0.007 \pm 0.004$. Let us underline that our three independent estimates of ϵ_M are fully consistent. Note also that this procedure is useful since the complete energy landscape is unknown and since the Maxwell point cannot be directly deduced from a single hysteresis diagram.

B. Reconstruction of the hysteresis loop at T = 0 K

Another important point is that we are able to recover the hysteresis loop at T = 0 K from the data recorded at $T \neq 0$. As shown by Fig. 11 (a), the energetic barriers are roughly linear functions of the confinement ϵ . We can thus extrapolate the data to estimate the values of $\epsilon_{up}(T = 0)$ and $\epsilon_{down}(T = 0)$ for which the energetic barriers $\Delta U_{L-B}(\epsilon)$ and $\Delta U_{B-L}(\epsilon)$ respectively vanish. These values correspond by definition to the thresholds and give the width of the hysteresis loop at T = 0 K. Moreover, the upper branch of the hysteresis loop is



FIG. 10: (Color online) (a) Plot of the mean residence time $\langle \tau_L \rangle$ (in s) for a line, as a function of temperature (in 10¹⁰ K), for $\epsilon = -0.01$. (b) Plot of the mean residence time $\langle \tau_B \rangle$ (in s) for a bubble, as a function of temperature (in 10¹⁰ K), for $\epsilon = -0.01$. The red dots are simulations data, the dashed blue line is the fit performed with Kramers' law.

quite accurately given by the position of the maximum of a typical PDF, since it varies slowly with temperature, while the lower branch is zero by definition of the line configuration. In Fig. 11 (b) we show an hysteresis loop at T = 0 K, reconstructed from measurements at $T \neq 0$, that compares rather well with the hysteresis loop evidenced by very low temperature simulations.



FIG. 11: (Color online) (a) Plot of the energetic barriers (10^{-12} J) as functions of the confinement parameter ϵ , for a transition from a line to a bubble $[\Delta U_{L-B}(\epsilon)]$, blue dots] and from a bubble to line $[\Delta U_{B-L}(\epsilon)]$, red squares]. (b) The dashed blue lines are determined from the extrapolation of the previous plots (see text for details) and indicate the width of the hysteresis loop. The shaded area is a measure of the uncertainty. For comparison the solid red line indicates the hysteresis loop obtained at very low temperature.

To conclude this discussion, let us emphasize an interesting opportunity given by our

analysis : We are able to obtain the local maximum $E_M(\epsilon)$ in the otherwise unreachable energy landscape. This is schematically illustrated in Fig. 12 (a). The energy $E_L(\epsilon)$ of a line and the energy $E_B(\epsilon)$ of a bubble may be calculated from the equilibrium configurations obtained in the simulations, and $E_M(\epsilon)$ is the energy amount that has to be spent in order to jump between the line and bubble configuration. Therefore, we may write

$$E_M(\epsilon) = E_L(\epsilon) + \Delta U_{L-B}(\epsilon) = E_B(\epsilon) + \Delta U_{B-L}(\epsilon).$$
(5)

We plot in Fig. 12 (b) these two sums. They are in very good agreement, especially if we consider the small values of the energetic barriers with respect to the configurations energies, which is of the order of 15 nJ. This result supports our analysis of the residence times as Kramer's times. Therefore, we propose this explicit procedure to determine the main features of the energy landscape in such many body systems.



FIG. 12: (Color online) (a) Schematic representation at a given ϵ of the energy landscape, where the arrows indicate the energetic barriers ΔU_{L-B} and ΔU_{B-L} . (b) Energy E_M (in 10⁻³ nJ) of the local maximum as a function of ϵ , see Eqn. (5). The blue dots display the reconstruction from the measurement of the energetic barrier ΔU_{L-B} and the energy of a line, and the red squares the reconstruction from the measurement of the energetic barrier ΔU_{B-L} and the energy of a bubble. The origin of the energy is the line configuration energy, E_L .

V. CONCLUSION

The subcritical pitchfork bifurcation studied in this paper is the zigzag transition in systems of particles that are transversally confined, with short range (non Coulombic) interactions and periodic boundary conditions in the longitudinal direction. As a consequence of the subcriticality, there is a finite range of transverse confinement stiffnesses for which the stable configuration at zero temperature is a bubble, where a localized zigzag pattern takes place in a finite part of the periodic cell. We have shown that these configurations remains stable for a wide temperature range.

At zero temperature, the transition between a straight line, for which all particles are aligned along the bottom of the transverse confinement potential, and a bubble is hysteretic. We have studied the influence of the thermal noise on this hysteresis loop. Its description is more difficult than at T = 0 K since thermally activated jumps between the two configurations always occur and the system can never stay forever in a unique metastable state. But since an actual simulation last a finite time, these jumps need to be taken into account only if their characteristic Kramers' times, $\tau_{L-B}(\epsilon, T)$ and $\tau_{B-L}(\epsilon, T)$ are smaller than the observation time τ_{obs} . The Kramers' times being temperature-dependent, two different regimes have to be considered according to the temperature value with respect to a critical temperature $T_c(\tau_{obs})$ such that $\tau_{L-B}(\epsilon, T_c) \sim \tau_{B-L}(\epsilon, T_c) \sim \tau_{obs}$. For $T < T_c(\tau_{obs})$ the duration of a simulation is short enough for the system to keep the memory of its initial condition. An hysteresis loop is still observed, with a width that decreases as the temperature increases toward $T_c(\tau_{obs})$. For $T > T_c(\tau_{obs})$ the memory of the initial condition is lost by stochastic jumps between the linear and bubble configurations evidenced on the duration τ_{obs} , and the hysteresis loop is no longer seen.

Moreover the study of the mean residence times in each configurations, measured in this intermittent regime, gives a unique opportunity to precisely determine the barrier height of the energy landscape of the considered system. Lastly from these values obtained at high temperature, we are able to determine the extension of the hysteresis loop which would exist at T = 0 K.

Appendix A: Numerics

The details on the numerical algorithm may be found in Ref. [35]. A short summary is presented here. We simulate identical point particles of mass m moving on a plane, submitted to a thermal bath at temperature T. The thermal bath is accounted for by a damping constant γ , and by random forces applied on each particle, with the statistical properties of uncorrelated white Gaussian noise. The particles are transversally confined in a quasi-1D geometry by a harmonic potential of stiffness β , whereas periodic boundary conditions are applied in the longitudinal direction. The particles interact with a repulsive potential, which in all simulations is a screened electrostatic interaction $U(r) = U_0 K_0(r/\lambda_0)$ with energy scale U_0 , with a characteristic length λ_0 and where K_0 is the modified Bessel function of index 0. The dynamics of the system is then simulated by the numerical integration of coupled Langevin equations [35].

Let L be the length of the periodic simulation cell, and let N be the number of particles. For all the simulations considered here, we have taken N = 128 and L = 240 mm, thus an inter particles distance d = 1.875 mm. The damping constant is $\gamma = 1$ s⁻¹. For such parameters the equilibrium properties of the system have been described by a subcritical pitchfork bifurcation (see Appendix B and Ref. [24]).

The focus in this paper is on the influence of the temperature on the subcritical bifurcation. Thus the range of temperature considered here is between 10^9 to 2. 10^{11} K, to be compared with the typical interaction energy between particles at our concerned density $U_{inter} \approx 0.117$ nJ ($U_{inter} \approx 8.46 \ 10^{12}$ K).

In some simulations, we vary the stiffness by steps. The typical duration of a stiffness step is 2. 10^5 time steps, and with a relative stiffness variation of 10^{-3} . To get enough statistics for the PDF of $h_m(\epsilon, T, t)$ we have also undertaken longer simulations (up to 10^7 time steps) at fixed confinement. Therefore, the analysis of the residence times has been obtained by the measure of at least thousand jumps between each configurations for every confinement studied. The smoothness of the resulting distributions (see Fig. 9) attest the quality of these measurements.

Appendix B: Bifurcation analysis of the zigzag transition.

The normal form for the zigzag transition has been established in Ref. [24]. In this appendix we summarize the main results.

1. Thermodynamic limit

The zigzag transition happens because it is energetically favorable for the particles to organize in a staggered row when the density 1/d increases or when the transverse stiffness

 β decreases. In this zigzag pattern, the particles alternate with a transverse coordinate $\pm h$ so that the nearest neighbor distance increases. Therefore, a decrease in the interaction energy eventually compensates the energy increase due to the transverse confining potential. Since the two symmetric zigzag patterns have the same energy, the zigzag transition is a pitchfork bifurcation which takes place at a critical value β_{ZZ} of the transverse stiffness. The physical origin of the normal form for this bifurcation is the coupling between the soft mode at the transition and the Goldstone mode due to the translational invariance of the infinite system.

A consistent expansion up to the power $\epsilon^{5/3}$, where ϵ is the distance to threshold $\beta = \beta_{ZZ}(1-\epsilon)$ gives the normal form as

$$m\ddot{h} = \beta_{ZZ}\epsilon h + g_3h^3 - g_5h^5 + g_1\partial^2 h/\partial x^2, \tag{B1}$$

where all distances are in units of d and where \dot{h} denote the time derivative of h. Knowing the interaction potential $U(r/\lambda_0)$ (see Appendix A), and restricting the interactions up to the second neighbors only, we have given in Ref. [24] the analytic expressions of β_{ZZ} and the coefficients g_i which are known functions of d. For $d/\lambda_0 > 2.04$ (see [24]) the coefficients g_3 and g_5 in Eqn. (B1) are positive and the pitchfork bifurcation is *subcritical*.

The line h = 0 is a stable equilibrium pattern when $\epsilon \leq 0$. The homogeneous zigzag pattern is stable when $-g_3^2/(4\beta_{ZZ}g_5) \leq \epsilon$, and its amplitude is

$$h_{ZZ}^2 = \frac{1}{2g_5} \left(g_3 + \sqrt{g_3^2 + 4\epsilon\beta_{ZZ}g_5} \right).$$
(B2)

In [24], we have shown that the normal form Eqn. (B1) allows inhomogeneous stationary solutions when $-3g_3^2/(16\beta_{ZZ}g_5) \leq \epsilon \leq 0$, such that

$$h(x) = \frac{h_{b-} h_{b+}}{\sqrt{(h_{b+}^2 - h_{b-}^2) \cosh^2\left(\sqrt{\frac{-\epsilon\beta_{ZZ}}{g_1}}x\right) + h_{b-}^2}},$$
(B3)

where

$$h_{b\pm}^2 = \frac{3}{2g_5} \left(\frac{g_3}{2} \pm \sqrt{\frac{g_3^2}{4} + \frac{4}{3}\epsilon g_5 \beta_{ZZ}} \right).$$
(B4)

We call these excitations *bubbles*. In the context of phase transitions, these localized excitations exhibit a phase coexistence between the two phases (aligned and zigzag), hence a first order transition. Their observation in our simulations evidences the subcriticallity of the zigzag bifurcation.

2. Periodic boundary conditions

The previous analysis may be extended to finite systems with periodic boundary conditions, because the periodicity implies a rotational invariance that induces a Goldstone mode. The normal form (B1) is thus relevant to periodic systems, assuming as for systems in the thermodynamic limit the vicinity of the bifurcation threshold ($|\epsilon| \ll 1$) and slow spatial variations ($|\partial h/\partial x| \ll h/d$).

Periodic boundary conditions are relevant for the simulations, which exhibit localized inhomogeneous patterns, as shown in [24] and in Sec. II. In Ref. [24], the simulations are done at very low temperature. The thermal fluctuations are very small, so that it is easy to measure the inter particle distance in that part of the inhomogeneous pattern where the particles are aligned (outside the bubble), which is found to be $d_{\text{eff}} > d$. It is therefore energetically favorable to have a localized excitation with nonzero transverse displacement h(x), because the particles that stay aligned with h = 0 are separated by a distance d_{eff} which decreases their interaction energy.

In the simulations, the inhomogeneous patterns are observed when the stiffness β_{simu} of the confining potential is such that $\beta_{\text{simu}} < \beta_{ZZ}(d)$. This defines a positive parameter ϵ_{simu} as $\beta_{\text{simu}} = \beta_{ZZ}(d)(1 - \epsilon_{\text{simu}})$. From the inter particle distance $d_{\text{eff}} > d$, we define an *effective* bifurcation parameter ϵ_{eff}

$$\epsilon_{\rm eff} = 1 - \frac{\beta_{\rm simu}}{\beta_{ZZ}(d_{\rm eff})}.$$
(B5)

We have shown in [24] that a bubble observed in a finite periodic system is very well described as a bubble in an infinite system, with an inter particle distance d_{eff} and a distance to threshold ϵ_{eff} .

At high temperature, it is less easy to measure d_{eff} because of the thermal fluctuations. The coupling between the two soft modes which is the basic mechanism responsible for the subcriticallity of the zigzag transition, thus for the possibility to observe the zigzag bubbles, lead to an analytic expression for the length of a periodic system of N particles that includes a bubble. This length has to be equal to the spatial period L, which is expressed by [24]

$$L = d_{\text{eff}} \left[N - \frac{2\alpha h_{b-} h_{b+}}{\kappa'} \operatorname{arctanh} \left(\frac{h_{b-}}{h_{b+}} \right) \right].$$
(B6)

Here α is a known function of d_{eff} which quantifies the mode coupling, and for a given value of the transverse stiffness β_{simu} this is an equation of the only unknown d_{eff} .



FIG. 13: (a) Plot of the numerical solution d_{eff} (mm) of (B6) as a function of ϵ_{simu} for N = 128and L = 240 mm. (b) Plot of the effective distance to threshold ϵ_{eff} as a function of ϵ_{simu} for N = 128 and L = 240 mm.

The numerical roots d_{eff} (mm) of (B6) and the corresponding effective distance to threshold ϵ_{eff} are plotted in Fig. 13 for several values of ϵ_{simu} [36]. We see that the calculated values of d_{eff} are always greater than d, and that ϵ_{eff} is negative, which is consistent with our description of the bubbles.

In Fig. 2 we compare the bubble patterns observed at various temperatures with the analytic expression (B3), calculated for the value of d_{eff} that is the root of (B6) for the value ϵ_{simu} and the particles number N used in the simulations. We see that when the temperature is low enough to avoid the pattern destruction by the thermal fluctuations the agreement is excellent.

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